Connecting via Winsock to STN

Welcome to STN International! Enter x:x

LOGINID:ssspta1619lxw

PASSWORD:

TERMINAL (ENTER 1, 2, 3, OR ?):2

```
Welcome to STN International
NEWS
                 Web Page URLs for STN Seminar Schedule - N. America
NEWS
                 "Ask CAS" for self-help around the clock
NEWS
         SEP 09
                 CA/CAplus records now contain indexing from 1907 to the
                 present.
NEWS
         AUG 05
                 New pricing for EUROPATFULL and PCTFULL effective
                 August 1, 2003
NEWS
         AUG 13
                 Field Availability (/FA) field enhanced in BEILSTEIN
                 Data available for download as a PDF in RDISCLOSURE
NEWS
        AUG 18
NEWS
      7
         AUG 18
                 Simultaneous left and right truncation added to PASCAL
NEWS
      8
         AUG 18
                 FROSTI and KOSMET enhanced with Simultaneous Left and Righ
                 Truncation
NEWS
     9
         AUG 18
                 Simultaneous left and right truncation added to ANABSTR
NEWS 10
         SEP 22
                 DIPPR file reloaded
NEWS 11
         SEP 25
                 INPADOC: Legal Status data to be reloaded
NEWS 12
        SEP 29
                 DISSABS now available on STN
NEWS 13
        OCT 10
                 PCTFULL: Two new display fields added
              OCTOBER 01 CURRENT WINDOWS VERSION IS V6.01a, CURRENT
NEWS EXPRESS
              MACINTOSH VERSION IS V6.0b(ENG) AND V6.0Jb(JP),
              AND CURRENT DISCOVER FILE IS DATED 23 SEPTEMBER 2003
NEWS HOURS
              STN Operating Hours Plus Help Desk Availability
NEWS INTER
              General Internet Information
NEWS LOGIN
              Welcome Banner and News Items
              Direct Dial and Telecommunication Network Access to STN
NEWS PHONE
NEWS WWW
              CAS World Wide Web Site (general information)
```

Enter NEWS followed by the item number or name to see news on that specific topic.

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FILE 'HOME' ENTERED AT 15:26:10 ON 15 OCT 2003

=> fil reg COST IN U.S. DOLLARS

FULL ESTIMATED COST

SINCE FILE TOTAL ENTRY SESSION 1.89 1.89

FILE 'REGISTRY' ENTERED AT 15:31:31 ON 15 OCT 2003 USE·IS SUBJECT TO THE TERMS OF YOUR STN CUSTOMER AGREEMENT. PLEASE SEE "HELP USAGETERMS" FOR DETAILS.

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STRUCTURE FILE UPDATES: 14 OCT 2003 HIGHEST RN 604736-26-7 DICTIONARY FILE UPDATES: 14 OCT 2003 HIGHEST RN 604736-26-7

TSCA INFORMATION NOW CURRENT THROUGH JULY 14, 2003

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```
=> e chlorocresol/cn
                   CHLOROCOPPER (1+)/CN
E1
             1
E2
             1
                   CHLOROCORONENE/CN
E3
             2 --> CHLOROCRESOL/CN
E4
             1
                   CHLOROCRESOL GREEN/CN
                   CHLOROCRESYL PHOSPHATE/CN
E5
             2
                   CHLOROCRIPTINE/CN
E6
             1
                   CHLOROCROTYL METHACRYLATE/CN
E7
             1
                   CHLOROCROTYLTUNGSTEN/CN
E8
             1
E9
             1
                   CHLOROCRUORIN (EUDISTYLIA VANCOUVERI A1 CHAIN PRECURSOR)/CN
                   CHLOROCRUORIN (SABELLASTARTE INDICA E-CHAIN PRECURSOR REDUCE
E10
             1
                   D)/CN
             1
                   CHLOROCRUOROHEME/CN
E11
                   CHLOROCRUOROHEMIN/CN
E12
=> s e3
             2 CHLOROCRESOL/CN
L1
=> d
     ANSWER 1 OF 2 REGISTRY COPYRIGHT 2003 ACS on STN
L1
     1321-10-4 REGISTRY
RN
CN
     Phenol, chloromethyl- (9CI)
                                   (CA INDEX NAME)
OTHER CA INDEX NAMES:
     Cresol, chloro- (7CI, 8CI)
CN
OTHER NAMES:
CN
     Chlorocresol
CN
     Chloromethylphenol
CN
     Monochlorocresol
DR
     86006-41-9, 29468-35-7, 31308-59-5
MF
     C7 H7 Cl O
     IDS, COM
CI
T.C
     STN Files:
                  ADISNEWS, AGRICOLA, BIOBUSINESS, BIOSIS, BIOTECHNO, CA,
       CAOLD, CAPLUS, CHEMLIST, CIN, EMBASE, PROMT, TOXCENTER, USPAT2,
       USPATFULL
                      EINECS**
     Other Sources:
         (**Enter CHEMLIST File for up-to-date regulatory information)
```



D1-Cl

D1-OH

D1-Me

```
=> d 2
     ANSWER 2 OF 2 REGISTRY COPYRIGHT 2003 ACS on STN
L1
     59-50-7 REGISTRY
RN
     Phenol, 4-chloro-3-methyl- (9CI) (CA INDEX NAME)
OTHER CA INDEX NAMES:
    m-Cresol, 4-chloro- (8CI)
OTHER NAMES:
CN
     1-Chloro-2-methyl-4-hydroxybenzene
CN
     2-Chloro-5-hydroxytoluene
CN
     3-Methyl-4-chlorophenol
CN
     4-Chloro-3-cresol
CN
     4-Chloro-3-methylphenol
     4-Chloro-5-methylphenol
CN
CN
     4-Chloro-m-cresol
CN
     6-Chloro-3-hydroxytoluene
CN
     Aptal
CN
     Baktol
CN
     Baktolan
CN
     Candaseptic
CN
     Chlorocresol
CN
     Neopredisan
CN
     NSC 4166
CN
     Ottafact
CN
     p-Chloro-m-cresol
CN
     para-Chloro-meta-cresol
CN
     Parol
CN
     PCMC
CN
     Peritonan
CN
     Preventol CMK
CN
     Raschit
CN
     Raschit K
CN
     Rasen-Anicon
FS
     3D CONCORD
MF
     C7 H7 C1 O
CI
     COM
LC
                  ADISNEWS, AGRICOLA, ANABSTR, AQUIRE, BEILSTEIN*, BIOBUSINESS,
     STN Files:
       BIOSIS, BIOTECHNO, CA, CANCERLIT, CAOLD, CAPLUS, CASREACT, CBNB,
       CHEMCATS, CHEMINFORMRX, CHEMLIST, CIN, CSCHEM, CSNB, DDFU, DRUGU,
       EMBASE, GMELIN*, HODOC*, HSDB*, IFICDB, IFIPAT, IFIUDB, IPA, MEDLINE,
       MRCK*, MSDS-OHS, NIOSHTIC, PHARMASEARCH, PIRA, PROMT, RTECS*, SPECINFO,
```

SYNTHLINE, TOXCENTER, ULIDAT, USAN, USPAT2, USPATFULL

178 REFERENCES IN FILE CA (1907 TO DATE)

179 REFERENCES IN FILE CAPLUS (1907 TO DATE) 4 REFERENCES IN FILE CAOLD (PRIOR TO 1967)

1 REFERENCES TO NON-SPECIFIC DERIVATIVES IN FILE CA

(*File contains numerically searchable property data)
Other Sources: DSL**, EINECS**, TSCA**, WHO
 (**Enter CHEMLIST File for up-to-date regulatory information)

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

1585 REFERENCES IN FILE CA (1907 TO DATE)
16 REFERENCES TO NON-SPECIFIC DERIVATIVES IN FILE CA
1587 REFERENCES IN FILE CAPLUS (1907 TO DATE)
22 REFERENCES IN FILE CAOLD (PRIOR TO 1967)

=> FIL REGISTRY

COST IN U.S. DOLLARS

SINCE FILE TOTAL ENTRY SESSION 7.98 9.87

FULL ESTIMATED COST

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STRUCTURE FILE UPDATES: 14 OCT 2003 HIGHEST RN 604736-26-7 DICTIONARY FILE UPDATES: 14 OCT 2003 HIGHEST RN 604736-26-7

TSCA INFORMATION NOW CURRENT THROUGH JULY 14, 2003

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=> S 59-50-7/RN

L2 1 59-50-7/RN

=> SET SMA OFF

SET COMMAND COMPLETED

=> DEL SEL Y

=> SEL RN

E1 THROUGH E1 ASSIGNED

=> SET SMA LOGIN

SET COMMAND COMPLETED

=> INDEX HSDB, HODOC, GMELIN

COST IN U.S. DOLLARS

SINCE FILE TOTAL ENTRY SESSION 0.48 10.35

FULL ESTIMATED COST

INDEX 'HSDB, HODOC, GMELIN' ENTERED AT 15:32:03 ON 15 OCT 2003

3 FILES IN THE FILE LIST IN STNINDEX

Enter SET DETAIL ON to see search term postings or to view search error messages that display as 0* with SET DETAIL OFF.

=> S E1 AND SLB/FA

1 FILE HSDB

1 FILE HODOC

2 FILES HAVE ONE OR MORE ANSWERS, 3 FILES SEARCHED IN STNINDEX

L3 OUE 59-50-7/BI AND SLB/FA

=> D RANK

F1 1 HSDB F2 1 HODOC

=> FIL HITS

COST IN U.S. DOLLARS

SINCE FILE TOTAL ENTRY SESSION

10.90

0.55

FULL ESTIMATED COST

FILE 'HSDB' ENTERED AT 15:32:08 ON 15 OCT 2003 COPYRIGHT (C) 2003 NATIONAL LIBRARY OF MEDICINE

FILE 'HODOC' ENTERED AT 15:32:08 ON 15 OCT 2003 COPYRIGHT (C) 2003 Crc Press, Inc. (CRC)

=> S L3

L4 2 L3

=> SET NOTICE 1 DISPLAY

NOTICE SET TO 1 U.S. DOLLAR FOR DISPLAY COMMAND SET COMMAND COMPLETED

=> D RN CN SLB 1-

YOU HAVE REQUESTED DATA FROM 2 ANSWERS - CONTINUE? Y/(N):Y THE ESTIMATED COST FOR THIS REQUEST IS 2.81 U.S. DOLLARS DO YOU WANT TO CONTINUE WITH THIS REQUEST? (Y)/N:Y

L4 ANSWER 1 OF 2 HSDB COPYRIGHT 2003 NLM on STN

```
59-50-7 HSDB
CAS Registry No. (RN):
Chemical Name (CN):
                              3-METHYL-4-CHLOROPHENOL
Synonyms (CN):
                              AI3-00075 **PEER REVIEWED**; APTAL **PEER
    REVIEWED**; BAKTOL **PEER REVIEWED**; BAKTOLAN **PEER REVIEWED**;
    CANDASEPTIC **PEER REVIEWED**; Caswell No 185A **PEER REVIEWED**;
    P-CHLOR-M-CRESOL **PEER REVIEWED**; CHLOROCRESOL **PEER REVIEWED**;
    4-CHLORO-M-CRESOL **PEER REVIEWED**; 6-Chloro-m-cresol **PEER REVIEWED**;
    P-CHLOROCRESOL **PEER REVIEWED**; p-chloro-m-cresol **PEER REVIEWED**;
    Chlorocresolo **PEER REVIEWED**; Chlorocresolum (Latin) **PEER REVIEWED**;
    4-CHLORO-1-HYDROXY-3-METHYLBENZENE **PEER REVIEWED**; 2-CHLORO-
    HYDROXYTOLUENE **PEER REVIEWED**; 2-CHLORO-5-HYDROXYTOLUENE **PEER
    REVIEWED**; 4-Chloro-3-hydroxytoluene **PEER REVIEWED**;
    6-CHLORO-3-HYDROXYTOLUENE **PEER REVIEWED**; Chlorokresolum **PEER
    REVIEWED**; 4-chloro-3-methyl phenol **PEER REVIEWED**;
    4-CHLORO-5-METHYLPHENOL **PEER REVIEWED**; Clorocresol (Spanish) **PEER
    REVIEWED**; chloro-3-cresol **PEER REVIEWED**; m-Cresol,4-chloro **PEER
    REVIEWED**; EPA Pesticide Chemical Code 064206 **PEER REVIEWED**; OTTAFACT
    **PEER REVIEWED**; Parachlorometacresol **PEER REVIEWED**; PARMETOL **PEER
    REVIEWED**; PAROL **PEER REVIEWED**; PCMC **PEER REVIEWED**; PERITONAN
    **PEER REVIEWED**; PHENOL, 4-CHLORO-3-METHYL- **PEER REVIEWED**; PREVENTOL
    CMK **PEER REVIEWED**; RASCHIT **PEER REVIEWED**; RASCHIT K **PEER
    REVIEWED**; RASEN-ANICON **PEER REVIEWED**
Shipping Name/No. (CN):
                              UN 2020 Chlorophenols, solid; IMO 6.1
                              Chlorophenols, solid
    Physical and Chemical Properties
Solubility (SLB):
    SOL IN FATS & OILS, ORGANIC SOLVENTS **PEER REVIEWED** [Sax, N.I. and
     R.J. Lewis, Sr. (eds.). Hawley's Condensed Chemical Dictionary. 11th ed.
     New York: Van Nostrand Reinhold Co., 1987. 268]
    MORE SOL IN HOT THAN IN COLD WATER; FREELY SOL IN ALCOHOL, BENZENE,
     CHLOROFORM, ETHER, FIXED OILS, TERPENES, ACETONE, AQ ALKALINE SOLN,
     PETROLEUM ETHER **PEER REVIEWED**
                                         [The Merck Index. 10th ed. Rahway, New
     Jersey: Merck Co., Inc., 1983. 299]
    3850 MG/L WATER AT 20 DEG C **PEER REVIEWED** [Callahan, M.A., M.W.
     Slimak, N.W. Gabel, et al. Water-Related Environmental Fate of 129
     Priority Pollutants. Volume I. EPA-440/4 79-029a. Washington, DC: U.S.
     Environmental Protection Agency, December 1979.p. 92-2]
     ANSWER 2 OF 2 HODOC COPYRIGHT 2003 CRC on STN
     59-50-7 HODOC
RN
CN
     Phenol, 4-chloro-3-methyl-
CN
     m-Cresol, 4-chloro-
CN
     Aptal
CN
     Baktol
CN
     Baktolan
CN
     Candaseptic
CN
     Chlorocresol
CN
     p-Chloro-m-cresol
CN
     para-Chloro-meta-cresol
CN
     4-Chloro-m-cresol
CN
     4-Chloro-3-cresol
CN
     2-Chloro-5-hydroxytoluene
CN
     6-Chloro-3-hydroxytoluene
     {\tt 1-Chloro-2-methyl-4-hydroxybenzene}
CN
CN
     4-Chloro-3-methylphenol
CN
     4-Chloro-5-methylphenol
CN
     3-Methyl-4-chlorophenol
CN
     Ottafact
CN
     Parmetol
CN
     Parol
CN
     PCMC
CN
    Peritonan
```

CN Preventol CMK
CN Raschit
CN Raschit K
CN Rasen-Anicon
CN Toluene, 2-chloro-5-hydroxy

Solubility

Description Solvent SLB.TX SLB.SOL

Slightly soluble water; chloroform

Soluble | alcohol; ether; petroleum ether

=> SET NOTICE LOGIN DISPLAY

NOTICE SET TO OFF FOR DISPLAY COMMAND SET COMMAND COMPLETED

=>

Connecting via Winsock to STN

Welcome to STN International! Enter x:x

LOGINID:ssspta1619lxw

PASSWORD:

TERMINAL (ENTER 1, 2, 3, OR ?):2

```
Welcome to STN International
NEWS
                 Web Page URLs for STN Seminar Schedule - N. America
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                 "Ask CAS" for self-help around the clock
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                 August 1, 2003
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         AUG 13
                 Field Availability (/FA) field enhanced in BEILSTEIN
         AUG 18
                 Data available for download as a PDF in RDISCLOSURE
NEWS
                 Simultaneous left and right truncation added to PASCAL
NEWS
      7
         AUG 18
NEWS
     8
         AUG 18
                 FROSTI and KOSMET enhanced with Simultaneous Left and Righ
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NEWS
     9
         AUG 18
                 Simultaneous left and right truncation added to ANABSTR
NEWS 10
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         SEP 29
                 DISSABS now available on STN
NEWS 13
        OCT 10
                 PCTFULL: Two new display fields added
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              MACINTOSH VERSION IS V6.0b(ENG) AND V6.0Jb(JP),
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              Direct Dial and Telecommunication Network Access to STN
NEWS WWW
              CAS World Wide Web Site (general information)
```

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FILE 'HOME' ENTERED AT 09:58:15 ON 15 OCT 2003

=> fil reg COST IN U.S. DOLLARS

SINCE FILE TOTAL ENTRY SESSION 0.21 0.21

FULL ESTIMATED COST

FILE 'REGISTRY' ENTERED AT 09:58:18 ON 15 OCT 2003
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```
=> e chlorocresol/cn
                   CHLOROCOPPER (1+)/CN
E1
             1
E2
             1
                   CHLOROCORONENE/CN
E3
             2 --> CHLOROCRESOL/CN
E4
             1
                   CHLOROCRESOL GREEN/CN
                   CHLOROCRESYL PHOSPHATE/CN
             2
E5
             1
                   CHLOROCRIPTINE/CN
E6
E7
             1
                   CHLOROCROTYL METHACRYLATE/CN
             1
                   CHLOROCROTYLTUNGSTEN/CN
E8
                   CHLOROCRUORIN (EUDISTYLIA VANCOUVERI A1 CHAIN PRECURSOR)/CN
E9
             1
                   CHLOROCRUORIN (SABELLASTARTE INDICA E-CHAIN PRECURSOR REDUCE
             1
E10
                   D)/CN
                   CHLOROCRUOROHEME/CN
E11
                   CHLOROCRUOROHEMIN/CN
E12
=> s e3
             2 CHLOROCRESOL/CN
L1
=> d
     ANSWER 1 OF 2 REGISTRY COPYRIGHT 2003 ACS on STN
L1
RN
     1321-10-4 REGISTRY
     Phenol, chloromethyl- (9CI)
                                   (CA INDEX NAME)
OTHER CA INDEX NAMES:
     Cresol, chloro- (7CI, 8CI)
OTHER NAMES:
     Chlorocresol
CN
     Chloromethylphenol
CN
     Monochlorocresol
CN
DR
     86006-41-9, 29468-35-7, 31308-59-5
MF
     C7 H7 Cl O
CI
     IDS, COM
                  ADISNEWS, AGRICOLA, BIOBUSINESS, BIOSIS, BIOTECHNO, CA,
LC
     STN Files:
       CAOLD, CAPLUS, CHEMLIST, CIN, EMBASE, PROMT, TOXCENTER, USPAT2,
       USPATFULL
                      EINECS**
     Other Sources:
         (**Enter CHEMLIST File for up-to-date regulatory information)
```



D1-C1

D1-0H

D1-Me

```
=> d 2
L1
     ANSWER 2 OF 2 REGISTRY COPYRIGHT 2003 ACS on STN
     59-50-7 REGISTRY
RN
     Phenol, 4-chloro-3-methyl- (9CI) (CA INDEX NAME)
OTHER CA INDEX NAMES:
    m-Cresol, 4-chloro- (8CI)
OTHER NAMES:
     1-Chloro-2-methyl-4-hydroxybenzene
     2-Chloro-5-hydroxytoluene
CN
     3-Methyl-4-chlorophenol
CN
     4-Chloro-3-cresol
     4-Chloro-3-methylphenol
CN
CN
     4-Chloro-5-methylphenol
CN
     4-Chloro-m-cresol
CN
     6-Chloro-3-hydroxytoluene
CN
     Aptal
CN
     Baktol
CN
     Baktolan
CN
     Candaseptic
CN
     Chlorocresol
CN
     Neopredisan
CN
     NSC 4166
CN
     Ottafact
CN
     p-Chloro-m-cresol
CN
     para-Chloro-meta-cresol
CN
     Parol
CN
     PCMC
CN
     Peritonan
CN
     Preventol CMK
CN
     Raschit
CN
     Raschit K
CN
     Rasen-Anicon
FS
     3D CONCORD
MF
     C7 H7 C1 O
CI
     COM
LC
     STN Files:
                  ADISNEWS, AGRICOLA, ANABSTR, AQUIRE, BEILSTEIN*, BIOBUSINESS,
       BIOSIS, BIOTECHNO, CA, CANCERLIT, CAOLD, CAPLUS, CASREACT, CBNB,
       CHEMCATS, CHEMINFORMRX, CHEMLIST, CIN, CSCHEM, CSNB, DDFU, DRUGU,
       EMBASE, GMELIN*, HODOC*, HSDB*, IFICDB, IFIPAT, IFIUDB, IPA, MEDLINE,
       MRCK*, MSDS-OHS, NIOSHTIC, PHARMASEARCH, PIRA, PROMT, RTECS*, SPECINFO,
```

SYNTHLINE, TOXCENTER, ULIDAT, USAN, USPAT2, USPATFULL

178 REFERENCES IN FILE CA (1907 TO DATE)

179 REFERENCES IN FILE CAPLUS (1907 TO DATE) 4 REFERENCES IN FILE CAOLD (PRIOR TO 1967)

1 REFERENCES TO NON-SPECIFIC DERIVATIVES IN FILE CA

(*File contains numerically searchable property data)
Other Sources: DSL**, EINECS**, TSCA**, WHO
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16 REFERENCES TO NON-SPECIFIC DERIVATIVES IN FILE CA
1587 REFERENCES IN FILE CAPLUS (1907 TO DATE)
22 REFERENCES IN FILE CAOLD (PRIOR TO 1967)

```
e chlorbutanol/cn
=>
                   CHLORBUPHAM/CN
E1
             1
                   CHLORBUTAMIDE/CN
E2
             1
             1 --> CHLORBUTANOL/CN
E3
E4
             1
                   CHLORBUTIN/CN
                   CHLORBUTINPENICILLIN/CN
E5
             1
                   CHLORBUTOL/CN
E6
             1
E7
             1
                   CHLORCAM/CN
                   CHLORCAM, MIXT. CONTG./CN
E8
             1
E9
             1
                   CHLORCARAGARD/CN
                   CHLORCARVACROL/CN
E10
             1
E11
             1
                   CHLORCHINALDIN/CN
E12
                   CHLORCHINALDOL/CN
=> s e3
L2
             1 CHLORBUTANOL/CN
=> d
     ANSWER 1 OF 1 REGISTRY COPYRIGHT 2003 ACS on STN
L2
RN
     57-15-8 REGISTRY
     2-Propanol, 1,1,1-trichloro-2-methyl- (6CI, 8CI, 9CI). (CA INDEX NAME)
CN
OTHER NAMES:
     .beta.,.beta.,.beta.-Trichloro-tert-butyl alcohol
CN
CN
     1,1,1-Trichloro-2-methyl-2-propanol
CN
     1,1,1-Trichloro-tert-butyl alcohol
CN
     2,2,2-Trichloro-1,1-dimethylethanol
     2-(Trichloromethyl)-2-propanol
CN
     Acetochlorone
CN
     Acetonchloroform
CN
CN
     Acetone chloroform
CN
     Anhydrous chlorobutanol
     Chlorbutanol
CN
CN
     Chlorbutol
CN
     Chloreton
CN
     Chloretone
CN
     Chlorobutanol
CN
     Chlortran
CN
     Clortran
CN
     Coliquifilm
```

```
CN
     Dentalone
CN
     Khloreton
CN
     Methaform
CN
     NSC 44794
CN
     NSC 4596
     NSC 5208
CN
CN
     Sedaform
CN
     Trichloro-tert-butyl alcohol
CN
     Trichlorobutanol
FS
     3D CONCORD
     C4 H7 Cl3 O
MF
CI
     COM
LC
     STN Files:
                    ADISNEWS, AGRICOLA, ANABSTR, AQUIRE, BEILSTEIN*, BIOBUSINESS,
        BIOSIS, BIOTECHNO, CA, CANCERLIT, CAOLD, CAPLUS, CASREACT, CBNB, CEN,
       CHEMCATS, CHEMINFORMRX, CHEMLIST, CIN, CSCHEM, CSNB, DDFU, DETHERM*, DIOGENES, DRUGU, EMBASE, GMELIN*, HODOC*, HSDB*, IFICDB, IFIPAT, IFIUDB, IPA, MEDLINE, MRCK*, MSDS-OHS, NAPRALERT, NIOSHTIC, PIRA, PROMT, RTECS*,
        SPECINFO, SYNTHLINE, TOXCENTER, USAN, USPATFULL, VETU
          (*File contains numerically searchable property data)
     Other Sources: DSL**, EINECS**, TSCA**, WHO
          (**Enter CHEMLIST File for up-to-date regulatory information)
    OH
   -c-ccl<sub>3</sub>
   Me
**PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT**
               897 REFERENCES IN FILE CA (1907 TO DATE)
                 4 REFERENCES TO NON-SPECIFIC DERIVATIVES IN FILE CA
               898 REFERENCES IN FILE CAPLUS (1907 TO DATE)
                30 REFERENCES IN FILE CAOLD (PRIOR TO 1967)
=> e methyl nicotinate/cn
E1
              1
                     METHYL NICKEL PYROPHOSPHATE/CN
E2
              1
                     METHYL NICLATE/CN
E3
              1 --> METHYL NICOTINATE/CN
E4
                     METHYL NICOTINATE 1-OXIDE/CN
              1
E5
              1
                     METHYL NICOTINATE HYDROCHLORIDE/CN
E6
              1
                     METHYL NICOTINATE N-OXIDE/CN
E7
              1
                     METHYL NICOTINATE-CARBOXYL-14C/CN
E8
              1
                     METHYL NICOTINIMIDATE/CN
E9
              1
                     METHYL NICOTINOYL LACTATE/CN
E10
              1
                     METHYL NICOTINOYLACETATE/CN
E11
              1
                     METHYL NICOTINOYLMETHYL SULFOXIDE/CN
E12
                     METHYL NIDORELLAURINATE/CN
=> s e3
L3
              1 "METHYL NICOTINATE"/CN
=> d
     ANSWER 1 OF 1 REGISTRY COPYRIGHT 2003 ACS on STN
L3
RN
     93-60-7 REGISTRY
     3-Pyridinecarboxylic acid, methyl ester (9CI) (CA INDEX NAME)
OTHER CA INDEX NAMES:
     Nicotinic acid, methyl ester (6CI, 7CI, 8CI)
OTHER NAMES:
```

```
CN
     3-(Carbomethoxy)pyridine
CN
     3-(Methoxycarbonyl)pyridine
CN
     m-(Methoxycarbonyl)pyridine
CN
     Methyl 3-pyridinecarboxylate
CN
     Methyl nicotinate
CN
     Nicometh
     NSC 13126
CN
     NSC 403799
CN
FS
     3D CONCORD
DR
     123574-61-8
     C7 H7 N O2
MF
CI
     COM
LC
                   AGRICOLA, ANABSTR, BEILSTEIN*, BIOBUSINESS, BIOSIS,
     STN Files:
       BIOTECHNO, CA, CANCERLIT, CAOLD, CAPLUS, CASREACT, CHEMCATS,
       CHEMINFORMRX, CHEMLIST, CIN, CSCHEM, CSNB, DDFU, DETHERM*, DIOGENES,
       DRUGU, EMBASE, HODOC*, IFICDB, IFIPAT, IFIUDB, IPA, MEDLINE, MRCK*, MSDS-OHS, NAPRALERT, NIOSHTIC, PROMT, RTECS*, SPECINFO, SYNTHLINE,
       TOXCENTER, USPAT2, USPATFULL
          (*File contains numerically searchable property data)
     Other Sources: DSL**, EINECS**, TSCA**
          (**Enter CHEMLIST File for up-to-date regulatory information)
```

RN

486-12-4 REGISTRY

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

```
790 REFERENCES IN FILE CA (1907 TO DATE)
2 REFERENCES TO NON-SPECIFIC DERIVATIVES IN FILE CA
791 REFERENCES IN FILE CAPLUS (1907 TO DATE)
45 REFERENCES IN FILE CAOLD (PRIOR TO 1967)
```

=> e triprolidine/cn E1 1 TRIPROAMYLIN/CN E2 1 TRIPROLIDIN/CN E3 1 --> TRIPROLIDINE/CN E4 TRIPROLIDINE HYDROCHLORIDE/CN 1 E5 1 TRIPROLIDINE HYDROCHLORIDE MONOHYDRATE/CN E6 1 TRIPROLIDINE OXALATE/CN E7 1 TRIPROP/CN E8 1 TRIPROPANOLAMINE/CN E9 1 TRIPROPANOLAMINE TRIACRYLATE/CN E10 1 TRIPROPANOLAMINE TRIS(3-(3,5-DI-TERT-BUTYL-4-HYDROXYPHENYL)P ROPIONATE) (ESTER)/CN 2 E11 TRIPROPARGYL CYANURATE/CN E12 TRIPROPARGYL ISOCYANURATE/CN => s e3 1 TRIPROLIDINE/CN L4 =>dANSWER 1 OF 1 REGISTRY COPYRIGHT 2003 ACS on STN 1.4

```
Pyridine, 2-[(1E)-1-(4-methylphenyl)-3-(1-pyrrolidinyl)-1-propenyl]- (9CI)
CN
     (CA INDEX NAME)
OTHER CA INDEX NAMES:
     Pyridine, 2-[1-(4-methylphenyl)-3-(1-pyrrolidinyl)-1-propenyl]-, (E)-
     Pyridine, 2-[3-(1-pyrrolidinyl)-1-p-tolylpropenyl]-, (E)- (8CI)
CN
OTHER NAMES:
     trans-1-(2-Pyridyl)-3-pyrrolidino-1-p-tolylprop-1-ene
CN
     trans-1-(4-Methylphenyl)-1-(2-pyridyl)-3-pyrrolidinoprop-1-ene
CN
     trans-2-[3-(1-Pyrrolidinyl)-1-p-tolypropenyl]pyridine
CN
CN
     Triprolidin
CN
     Triprolidine
CN
     Tripyrolidine
FS
     STEREOSEARCH
MF
     C19 H22 N2
CI
     COM
LC
     STN Files:
                  ADISNEWS, AGRICOLA, ANABSTR, BEILSTEIN*, BIOBUSINESS, BIOSIS,
       BIOTECHNO, CA, CANCERLIT, CAOLD, CAPLUS, CHEMCATS, CHEMLIST, CIN,
       CSCHEM, DDFU, DIOGENES, DRUGU, EMBASE, HSDB*, IFICDB, IFIUDB, IPA,
       MEDLINE, MRCK*, PROMT, RTECS*, SPECINFO, TOXCENTER, USAN, USPATFULL
         (*File contains numerically searchable property data)
     Other Sources:
                     EINECS**, WHO
         (**Enter CHEMLIST File for up-to-date regulatory information)
```

Double bond geometry as shown.

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

347 REFERENCES IN FILE CA (1907 TO DATE)
6 REFERENCES TO NON-SPECIFIC DERIVATIVES IN FILE CA
348 REFERENCES IN FILE CAPLUS (1907 TO DATE)
3 REFERENCES IN FILE CAOLD (PRIOR TO 1967)

```
=> e promethazine/cn
                   PROMETHATE(5-), TETRAKIS(3,5-DINITRO-1,2-BENZENEDIOLATO(2-)-
E1
                   O1,O2)-, 9-(2-CARBOXYPHENYL)-3,6-BIS(DIETHYLAMINO)XANTHYLIUM
                    TETRAHYDROGEN/CN
                   PROMETHATE(5-), TETRAKIS(3,5-DINITRO-1,2-BENZENEDIOLATO(2-)-
E2
                   O1,O2)-, N-(9-(2-CARBOXYPHENYL)-6-(DIETHYLAMINO)-3H-XANTHEN-
                   3-YLIDENE) - N-ETHYLETHANAMINIUM TETRAHYDROGEN/CN
E3
             1 --> PROMETHAZINE/CN
                   PROMETHAZINE 3,4,5-TRIMETHOXYBENZOATE/CN
E4
             1
E5
             1
                   PROMETHAZINE 5,5-DIOXIDE/CN
E6
             .1
                   PROMETHAZINE 5-OXIDE/CN
E7
             1
                   PROMETHAZINE 5-SULFOXIDE/CN
                   PROMETHAZINE 8-CHLOROTHEOPHYLLINATE/CN
E8
```

```
PROMETHAZINE CHLORIDE/CN
E9
              1
                    PROMETHAZINE HYDROBROMIDE/CN
E10
              1
                    PROMETHAZINE HYDROCHLORIDE/CN
E11
              1
                    PROMETHAZINE MALEATE/CN
E12
              1
=> s e3
              1 PROMETHAZINE/CN
L5
=> d
     ANSWER 1 OF 1 REGISTRY COPYRIGHT 2003 ACS on STN
L5
RN
     60-87-7 REGISTRY
     10H-Phenothiazine-10-ethanamine, N,N,.alpha.-trimethyl- (9CI) (CA INDEX
CN
     NAME)
OTHER CA INDEX NAMES:
     Phenothiazine, 10-[2-(dimethylamino)propyl]- (8CI)
OTHER NAMES:
CN
     (.+-.)-Promethazine
     (2-Dimethylamino-2-methyl)ethyl-N-dibenzoparathiazine
CN
CN
     10-[2-(Dimethylamino)propyl]phenothiazine
CN
     Dimapp
CN
     Diphergan
CN
     NSC 30321
CN
     Proazamine
CN
     Procit
CN
     Prometazin
CN
     Promethazine
     Protazine
CN
CN
     Prothazin
CN
     RP 3277
     Vallergine
CN
     3D CONCORD
FS
     73745-50-3
DR
MF
     C17 H20 N2 S
CI
     COM
LC
     STN Files:
                   ADISNEWS, AGRICOLA, ANABSTR, BEILSTEIN*, BIOBUSINESS, BIOSIS,
       BIOTECHNO, CA, CABA, CANCERLIT, CAOLD, CAPLUS, CASREACT, CBNB, CHEMCATS, CHEMLIST, CIN, CSCHEM, DDFU, DIOGENES, DRUGU, EMBASE, GMELIN*, HODOC*,
       HSDB*, IFICDB, IFIPAT, IFIUDB, IPA, MEDLINE, MRCK*, NIOSHTIC,
       PHARMASEARCH, PROMT, RTECS*, SPECINFO, TOXCENTER, USAN, USPAT2,
       USPATFULL, VETU
          (*File contains numerically searchable property data)
                        EINECS**, WHO
     Other Sources:
          (**Enter CHEMLIST File for up-to-date regulatory information)
```

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

3322 REFERENCES IN FILE CA (1907 TO DATE)
48 REFERENCES TO NON-SPECIFIC DERIVATIVES IN FILE CA
3324 REFERENCES IN FILE CAPLUS (1907 TO DATE)

43 REFERENCES IN FILE CAOLD (PRIOR TO 1967)

```
=> e trimeprazine/cn
                   TRIMEPRANOL FUMARATE/CN
E1
             1
                   TRIMEPRANOL RETARD/CN
E2
             1
             1 --> TRIMEPRAZINE/CN
E3
                   TRIMEPRAZINE 5,5-DIOXIDE/CN
E4
             1
                   TRIMEPRAZINE HYDROCHLORIDE/CN
E5
             1
                   TRIMEPRAZINE MALEATE/CN
E6
             1
                   TRIMEPRAZINE RADICAL CATION/CN
E7
             1
E8
             1
                   TRIMEPRAZINE SULFOXIDE/CN
                   TRIMEPRAZINE TARTRATE/CN
E9
             1
E10
             1
                   TRIMEPRIMINE/CN
E11
             1
                   TRIMEPRIMINE HYDROCHLORIDE/CN
E12
             1
                   TRIMEPRIMINE MALEATE/CN
=> s e3
             1 TRIMEPRAZINE/CN
L6
=> d
     ANSWER 1 OF 1 REGISTRY COPYRIGHT 2003 ACS on STN
L6
     84-96-8 REGISTRY
RN
CN
     10H-Phenothiazine-10-propanamine, N,N,.beta.-trimethyl- (9CI) (CA INDEX
OTHER CA INDEX NAMES:
    Phenothiazine, 10-[3-(dimethylamino)-2-methylpropyl]- (6CI, 8CI)
OTHER NAMES:
CN
     (.+-.)-Alimemazine
CN
     (.+-.)-Trimeprazine
     10-(2-Methyl-3-dimethylaminopropyl)phenothiazine
CN
     10-[3-(Dimethylamino)-2-methylpropyl]phenothiazine
CN
CN
    Alimemazine
CN
    Alimezine
CN
    Bayer 1219
CN
     dl-Trimeprazine
CN
     Methylpromazine
CN
     Teralen
CN
     Teralene
     Trimeprazine
CN
FS
     3D CONCORD
DR
     35309-60-5, 47138-21-6
MF
     C18 H22 N2 S
CI
     COM
LC
     STN Files:
                  ADISNEWS, AGRICOLA, ANABSTR, BEILSTEIN*, BIOBUSINESS, BIOSIS,
       BIOTECHNO, CA, CANCERLIT, CAOLD, CAPLUS, CASREACT, CHEMLIST, CIN, DDFU,
       DRUGU, EMBASE, HSDB*, IPA, MEDLINE, MRCK*, NIOSHTIC, RTECS*, SPECINFO,
       TOXCENTER, USAN, USPATFULL, VETU
         (*File contains numerically searchable property data)
                     EINECS**; WHO
     Other Sources:
         (**Enter CHEMLIST File for up-to-date regulatory information)
          Me
Me2N-CH2-CH-CH2
```

```
548 REFERENCES IN FILE CAPLUS (1907 TO DATE)
              48 REFERENCES IN FILE CAOLD (PRIOR TO 1967)
=> s sulfiram/cn
             1 SULFIRAM/CN
=> s e3
L8
             1 TRIMEPRAZINE/CN
=> d 17
     ANSWER 1 OF 1 REGISTRY COPYRIGHT 2003 ACS on STN
     95-05-6 REGISTRY
RN
     Thiodicarbonic diamide ([(H2N)C(S)]2S), tetraethyl- (9CI) (CA INDEX NAME)
OTHER CA INDEX NAMES:
    Sulfide, bis(diethylthiocarbamoyl) (6CI, 7CI, 8CI)
OTHER NAMES:
    Bis (diethylthiocarbamoyl) sulfide
CN
CN
     Bis(diethylthiocarbamyl) sulfide
CN
     Bis (N, N-diethylthiocarbamoyl) sulfide
     Carbamodithioic acid, diethyl-, anhydrosulfide
CN
     Kutka
CN
CN
     Kutkasin
    Methanethioamide, 1,1'-thiobis[N,N-diethyl-
CN
CN
     Monosulfiram
     NSC 36731
CN
CN
     Sanigal
CN
     Sarcocide B
     Sulfide, bis[(diethylamino)thioxomethyl]
CN
CN
     Sulfiram
CN
     Sulfirame
CN
     Sulfiramum
CN
     Tetmos
CN
     Tetmosol
CN
     Tetraethylthiuram monosulfide
CN
     Tetrucid
CN
     TTMS
     3D CONCORD
FS
     C10 H20 N2 S3
MF
CI
LC
     STN Files:
                  ADISNEWS, AGRICOLA, AQUIRE, BEILSTEIN*, BIOSIS, CA, CABA,
       CAOLD, CAPLUS, CASREACT, CHEMCATS, CHEMLIST, CSCHEM, DDFU, DRUGU,
       EMBASE, HODOC*, IPA, MEDLINE, MRCK*, PHARMASEARCH, PROMT, RTECS*,
       TOXCENTER, USAN, USPATFULL
         (*File contains numerically searchable property data)
     Other Sources: EINECS**, WHO
         (**Enter CHEMLIST File for up-to-date regulatory information)
Et2N-C-S-C-NEt2
**PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT**
              76 REFERENCES IN FILE CA (1907 TO DATE)
              76 REFERENCES IN FILE CAPLUS (1907 TO DATE)
               8 REFERENCES IN FILE CAOLD (PRIOR TO 1967)
```

546 REFERENCES IN FILE CA (1907 TO DATE)

10 REFERENCES TO NON-SPECIFIC DERIVATIVES IN FILE CA

```
=> e oxybutynin/cn
E1
                   OXYBUPROCAINE/CN
             1
                   OXYBUPROCAINE HYDROCHLORIDE/CN
E2
             1
             1 --> OXYBUTYNIN/CN
E3
                   OXYBUTYNIN CHLORIDE/CN
E4
             1
                   OXYBUTYNIN HYDROCHLORIDE/CN
E5
             1
                   OXYBUTYNIN N-OXIDE/CN
E6
             1
E7
             1
                   OXYCADYSTIN/CN
                   OXYCAINE/CN
E8
             1
                   OXYCAINE 12/CN
E9
             1
                   OXYCAINE 4/CN
E10
             1
                   OXYCAINE HYDROCHLORIDE/CN
E11
             1
                   OXYCAINE PENICILLIN-G SALT/CN
E12
=> s e3
             1 OXYBUTYNIN/CN
L9
=>d
L9
     ANSWER 1 OF 1 REGISTRY COPYRIGHT 2003 ACS on STN
RN
     5633-20-5 REGISTRY
CN ·
     Benzeneacetic acid, .alpha.-cyclohexyl-.alpha.-hydroxy-,
     4-(diethylamino)-2-butynyl ester (9CI) (CA INDEX NAME)
OTHER CA INDEX NAMES:
CN
     2-Butyn-1-ol, 4-(diethylamino)-, .alpha.-phenylcyclohexaneglycolate
     (ester)
CN
     Cyclohexaneglycolic acid, .alpha.-phenyl-, 4-(diethylamino)-2-butynyl
     ester (8CI)
OTHER NAMES:
CN
     (.+-.)-Oxybutynin
CN
     (RS) -Oxybutynin
CN
     4-Diethylamino-2-butynyl .alpha.-phenylcyclohexaneglycolate
CN
     Ditropan
CN
     Oxybutynin
CN
     Oxytrol
FS
     3D CONCORD
DR
     119579-36-1
MF
     C22 H31 N O3
CI
     COM
LC
     STN Files:
                  ADISINSIGHT, ADISNEWS, ANABSTR, BEILSTEIN*, BIOBUSINESS,
       BIOSIS, BIOTECHNO, CA, CANCERLIT, CAPLUS, CBNB, CEN, CHEMCATS, CIN,
       CSCHEM, DDFU, DIOGENES, DRUGNL, DRUGPAT, DRUGU, EMBASE, HSDB*, IPA,
       MEDLINE, MRCK*, PHAR, PROMT, RTECS*, SPECINFO, TOXCENTER, USAN, USPAT2,
       USPATFULL
         (*File contains numerically searchable property data)
     Other Sources:
```

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

295 REFERENCES IN FILE CA (1907 TO DATE)

6 REFERENCES TO NON-SPECIFIC DERIVATIVES IN FILE CA 296 REFERENCES IN FILE CAPLUS (1907 TO DATE)

```
=> e tetosterone enanthate/cn
E1
             1
                   TETORON UL 9/CN
                   TETORON V/CN
E2
             1
E3
             0 --> TETOSTERONE ENANTHATE/CN
                   TETOXACIN/CN
E4
             1
             1
                   TETP/CN
E5
E6
             1
                   TETQ FAMILY GTPASE (PLASMODIUM FALCIPARUM STRAIN 3D7 GENE PF
                   L1710C)/CN
E7
                   TETR FAMILY HTH TRANSCRIPTIONAL REGULATOR (CLOSTRIDIUM ACETO
                   BUTYLICUM STRAIN ATCC 824 GENE CAP0046)/CN
                   TETR FAMILY REGULATORY PROTEIN (BORDETELLA BRONCHISEPTICA ST
E8
                   RAIN RB50 GENE BB0576)/CN
E9
                   TETR FAMILY REGULATORY PROTEIN (BORDETELLA PARAPERTUSSIS STR
                   AIN 12822 GENE BPP0570)/CN
                   TETR FAMILY REGULATORY PROTEIN (STREPTOMYCES COELICOLOR STRA
E10
                   IN A3(2) GENE 2SC10A7.12)/CN
                   TETR FAMILY TRANSCRIPTIONAL REGULATOR (BORDETELLA BRONCHISEP
E11
                   TICA STRAIN RB50 GENE BB0631)/CN
E12
                   TETR FAMILY TRANSCRIPTIONAL REGULATOR (BORDETELLA PARAPERTUS
                   SIS STRAIN 12822 GENE BPP0625)/CN
=> e testosterone enanthate/cn
E1
             1
                   TESTOSTERONE DIAZOACETATE/CN
E2
             1
                   TESTOSTERONE DIPROPIONATE/CN
E3
             1 --> TESTOSTERONE ENANTHATE/CN
E4
                   TESTOSTERONE ENANTHATE, HYDROXYPROGESTERONE CAPROATE, ESTRAD
                   IOL VALERATE/CN
                   TESTOSTERONE ENANTHATE-ESTRADIOL VALERATE MIXTURE/CN
E5
             1
E6
             1
                   TESTOSTERONE ENANTHATE-MEDROXYPROGESTERONE ACETATE MIXT./CN
E7
             1
                   TESTOSTERONE ENOL BIS (HEPTAFLUOROBUTYRATE) / CN
E8
             1
                   TESTOSTERONE ENOL BIS (TRIMETHYLSILYL) ETHER/CN
             1
                   TESTOSTERONE ENOL DIACETATE/CN
E9
                   TESTOSTERONE ETHYL DIAZOMALONATE/CN
E10
E11
             1
                   TESTOSTERONE ETHYLENE KETAL/CN
E12
                   TESTOSTERONE FLUOROACETATE/CN
=> s e3
L10
             1 "TESTOSTERONE ENANTHATE"/CN
=> d
L10
     ANSWER 1 OF 1 REGISTRY COPYRIGHT 2003 ACS on STN
RN
     315-37-7 REGISTRY
     Androst-4-en-3-one, 17-[(1-oxoheptyl)oxy]-, (17.beta.)- (9CI) (CA INDEX
     NAME)
OTHER CA INDEX NAMES:
    Testosterone, heptanoate (6CI, 8CI)
OTHER NAMES:
CN
     17.beta.-Enanthoxyandrost-4-en-3-one
     17.beta.-Hydroxyandrost-4-en-3-one enanthate
CN
CN
     4-Androsten-3-one 17.beta.-enanthate
CN
     Andro LA 200
CN
     Androtardyl
CN
     Delatestryl
CN
     Everone
CN
    NSC 17591
     Primoteston-Depot
CN
CN
     Reposo TMD
CN
     Testenate
CN
     Testinon
```

```
CN
     Testo-Enant
CN
     Testosterone 17-enanthate
CN
     Testosterone enanthate
CN
     Testosterone heptylate
CN
     Testosterone oenanthate
FS
     STEREOSEARCH
DR
     11111-10-7
MF
     C26 H40 O3
CI
     COM
                    ADISINSIGHT, ADISNEWS, AGRICOLA, ANABSTR, BEILSTEIN*,
LC
     STN Files:
        BIOBUSINESS, BIOSIS, BIOTECHNO, CA, CABA, CANCERLIT, CAOLD, CAPLUS,
       CBNB, CHEMCATS, CHEMLIST, CIN, CSCHEM, DDFU, DIOGENES, DRUGU, EMBASE, IFICDB, IFIPAT, IFIUDB, IPA, MEDLINE, MRCK*, MSDS-OHS, PROMT, RTECS*,
        TOXCENTER, USAN, USPAT2, USPATFULL, VETU
          (*File contains numerically searchable property data)
                        EINECS**
     Other Sources:
          (**Enter CHEMLIST File for up-to-date regulatory information)
```

Absolute stereochemistry. Rotation (+).

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

489 REFERENCES IN FILE CA (1907 TO DATE)
490 REFERENCES IN FILE CAPLUS (1907 TO DATE)
45 REFERENCES IN FILE CAOLD (PRIOR TO 1967)

```
=> e choline salicylate/cn
                   CHOLINE REINECKATE, PHENYL PHOSPHATE, ESTER WITH 1,2-DIHEXAN
E1
                   OIN/CN
E2
                   CHOLINE REINECKATE, PHENYL PROPYL PHOSPHATE/CN
             1
             1 --> CHOLINE SALICYLATE/CN
E3
E4
                   CHOLINE SALICYLIC ACID SALT/CN
             1
E5
                   CHOLINE SALT WITH PHENOLPHTHALIN/CN
             1
E6
                   CHOLINE SENECIOATE, 2,2',4,4',6,6'-HEXANITRODIPHENYLAMINE DE
             1
                   RIV./CN
E7
             7
                   CHOLINE SENNOSIDE A/CN
                   CHOLINE SENNOSIDE B/CN
E8
             1
E9
                   CHOLINE SUBSALICYLATE/CN
             1
E10
                   CHOLINE SUCCINATE DICHLORIDE/CN
             1
E11
                   CHOLINE SUCCINATE DICHLORIDE DIHYDRATE/CN
             1
                   CHOLINE SULFADIMERAZINE-N-1-PROPANESULFONATE/CN
E12
=> s e3
             1 "CHOLINE SALICYLATE"/CN
L11
```

```
ANSWER 1 OF 1 REGISTRY COPYRIGHT 2003 ACS on STN
L11
     2016-36-6 REGISTRY
RN
     Ethanaminium, 2-hydroxy-N,N,N-trimethyl-, salt with 2-hydroxybenzoic acid
CN
     (1:1) (9CI) (CA INDEX NAME)
OTHER CA INDEX NAMES:
     Benzoic acid, 2-hydroxy-, ion(1-), 2-hydroxy-N,N,N-trimethylethanaminium
     (9CI)
CN
     Choline salicylate (6CI)
     Choline, salicylate (salt) (7CI, 8CI)
CN
CN
     Salicylic acid, ion(1-), choline (8CI)
OTHER NAMES:
CN
     (2-Hydroxyethyl) trimethylammonium salicylate
CN
     Actasal
CN
     Arret
CN
     Arthropan
CN
     Artrobione
CN
     Audax
CN
     Choline salicylic acid salt
CN
     Choline subsalicylate
CN
     Mundisal
CN
     Salicol
CN
     Salicylic acid choline salt
CN
     Satibon
CN
     Syrap
DR
     54391-51-4
     C7 H5 O3 . C5 H14 N O
MF
CI
       TN Files: ADISNEWS, AGRICOLA, ANABSTR, BEILSTEIN*, BIOBUSINESS, BIOSIS, BIOTECHNO, CA, CANCERLIT, CAOLD, CAPLUS, CASREACT, CBNB, CHEMLIST, CIN,
LC
     STN Files:
       CSCHEM, DDFU, DIOGENES, DRUGU, EMBASE, IFICDB, IFIPAT, IFIUDB, IPA,
       MEDLINE, MRCK*, PHARMASEARCH, PROMT, RTECS*, TOXCENTER, USAN, USPATFULL
          (*File contains numerically searchable property data)
                      EINECS**, NDSL**, TSCA**, WHO
     Other Sources:
          (**Enter CHEMLIST File for up-to-date regulatory information)
     CM
          1
     CRN 63-36-5
     CMF C7 H5 O3
       CO2 -
       ОН
     CM
          2
     CRN 62-49-7
     CMF C5 H14 N O
Me_3+N-CH_2-CH_2-OH
              113 REFERENCES IN FILE CA (1907 TO DATE)
                8 REFERENCES TO NON-SPECIFIC DERIVATIVES IN FILE CA
              114 REFERENCES IN FILE CAPLUS (1907 TO DATE)
               14 REFERENCES IN FILE CAOLD (PRIOR TO 1967)
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